



# Full wwPDB X-ray Structure Validation Report ⓘ

Feb 1, 2016 – 07:20 PM GMT

PDB ID : 4OJZ  
Title : Crystal Structure of Alg17c Mutant Y258A Complexed with Alginate Trisaccharide  
Authors : Nair, S.K.; Park, D.S.  
Deposited on : 2014-01-21  
Resolution : 1.90 Å(reported)

This is a Full wwPDB X-ray Structure Validation Report for a publicly released PDB entry.  
We welcome your comments at [validation@mail.wwpdb.org](mailto:validation@mail.wwpdb.org)  
A user guide is available at  
<http://wwpdb.org/validation/2016/XrayValidationReportHelp>  
with specific help available everywhere you see the ⓘ symbol.

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The following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

MolProbity : 4.02b-467  
Mogul : 1.7 (RC4), CSD as536be (2015)  
Xtriage (Phenix) : 1.9-1692  
EDS : rb-20026688  
Percentile statistics : 20151230.v01 (using entries in the PDB archive December 30th 2015)  
Refmac : 5.8.0135  
CCP4 : 6.5.0  
Ideal geometry (proteins) : Engh & Huber (2001)  
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)  
Validation Pipeline (wwPDB-VP) : trunk26865

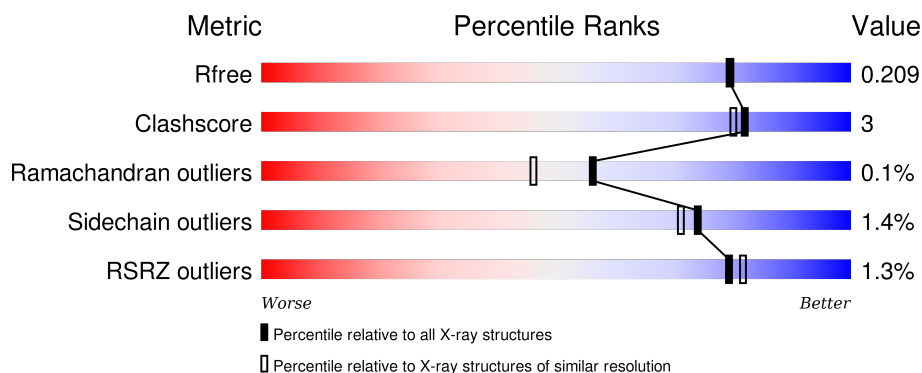
# 1 Overall quality at a glance ⓘ

The following experimental techniques were used to determine the structure:

## *X-RAY DIFFRACTION*

The reported resolution of this entry is 1.90 Å.

Percentile scores (ranging between 0-100) for global validation metrics of the entry are shown in the following graphic. The table shows the number of entries on which the scores are based.



Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
$R_{free}$	91344	4755 (1.90-1.90)
Clashscore	102246	5398 (1.90-1.90)
Ramachandran outliers	100387	5338 (1.90-1.90)
Sidechain outliers	100360	5339 (1.90-1.90)
RSRZ outliers	91569	4766 (1.90-1.90)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments on the lower bar indicate the fraction of residues that contain outliers for  $\geq 3$ , 2, 1 and 0 types of geometric quality criteria. A grey segment represents the fraction of residues that are not modelled. The numeric value for each fraction is indicated below the corresponding segment, with a dot representing fractions  $\leq 5\%$ . The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density. The numeric value is given above the bar.

Mol	Chain	Length	Quality of chain
1	A	736	<div> <div>%</div> <div> <div></div> <div>90%</div> <div>6%</div> <div>.</div> </div> </div>
1	B	736	<div> <div>2%</div> <div> <div></div> <div>89%</div> <div>7%</div> <div>.</div> </div> </div>

The following table lists non-polymeric compounds, carbohydrate monomers and non-standard residues in protein, DNA, RNA chains that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
3	MAV	A	803	X	-	-	-
3	MAV	B	803	X	-	-	-
4	EDO	B	805	-	-	-	X

## 2 Entry composition

There are 5 unique types of molecules in this entry. The entry contains 12602 atoms, of which 0 are hydrogens and 0 are deuteriums.

In the tables below, the ZeroOcc column contains the number of atoms modelled with zero occupancy, the AltConf column contains the number of residues with at least one atom in alternate conformation and the Trace column contains the number of residues modelled with at most 2 atoms.

- Molecule 1 is a protein called Putative alginate lyase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	706	Total	C	N	O	S	0	0	0
			5549	3531	938	1067	13			
1	B	705	Total	C	N	O	S	0	0	0
			5538	3525	934	1066	13			

There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	258	ALA	TYR	ENGINEERED MUTATION	UNP Q21FJ0
B	258	ALA	TYR	ENGINEERED MUTATION	UNP Q21FJ0

- Molecule 2 is ZINC ION (three-letter code: ZN) (formula: Zn).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	B	1	Total	Zn	0	0
			1	1		
2	A	1	Total	Zn	0	0
			1	1		

- Molecule 3 is a polymer of unknown type called SUGAR (3-MER).

Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	A	3	Total	C	O	0	0
			36	18	18		
3	B	3	Total	C	O	0	0
			36	18	18		

- Molecule 4 is 1,2-ETHANEDIOL (three-letter code: EDO) (formula: C<sub>2</sub>H<sub>6</sub>O<sub>2</sub>).



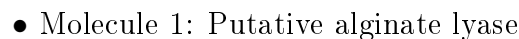
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
4	A	1	Total	C	O	0	0
			4	2	2		
4	B	1	Total	C	O	0	0
			4	2	2		

- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	731	Total	O	0	0
			731	731		
5	B	702	Total	O	0	0
			702	702		



- Molecule 1: Putative alginate lyase



## 4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	85.82Å 128.62Å 88.13Å 90.00° 111.81° 90.00°	Depositor
Resolution (Å)	24.00 – 1.90 38.85 – 1.90	Depositor EDS
% Data completeness (in resolution range)	99.6 (24.00-1.90) 99.6 (38.85-1.90)	Depositor EDS
$R_{merge}$	(Not available)	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	2.49 (at 1.89Å)	Xtriage
Refinement program	REFMAC 5.7.0029	Depositor
R, $R_{free}$	0.179 , 0.209 0.179 , 0.209	Depositor DCC
$R_{free}$ test set	6939 reflections (5.26%)	DCC
Wilson B-factor (Å <sup>2</sup> )	18.5	Xtriage
Anisotropy	0.689	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.37 , 50.6	EDS
Estimated twinning fraction	0.016 for l,-k,h	Xtriage
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.49$ , $\langle L^2 \rangle = 0.32$	Xtriage
Outliers	0 of 138807 reflections	Xtriage
$F_o, F_c$ correlation	0.95	EDS
Total number of atoms	12602	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	22.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.46% of the height of the origin peak. No significant pseudotranslation is detected.*

<sup>1</sup> Intensities estimated from amplitudes.

<sup>2</sup> Theoretical values of  $\langle |L| \rangle$ ,  $\langle L^2 \rangle$  for acentric reflections are 0.5, 0.375 respectively for untwinned datasets, and 0.333, 0.2 for perfectly twinned datasets.

## 5 Model quality

### 5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: LGU, ZN, BEM, EDO, MAV

The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with  $|Z| > 5$  is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	# Z  >5	RMSZ	# Z  >5
1	A	0.34	0/5677	0.53	1/7705 (0.0%)
1	B	0.34	0/5666	0.51	0/7691
All	All	0.34	0/11343	0.52	1/15396 (0.0%)

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand. A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
3	A	1	0
3	B	1	0
All	All	2	0

There are no bond length outliers.

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	112	LEU	CA-CB-CG	5.03	126.86	115.30

All (2) chirality outliers are listed below:

Mol	Chain	Res	Type	Atom
3	A	803	MAV	C1
3	B	803	MAV	C1

There are no planarity outliers.



## 5.2 Too-close contacts ⓘ

In the following table, the Non-H and H(model) columns list the number of non-hydrogen atoms and hydrogen atoms in the chain respectively. The H(added) column lists the number of hydrogen atoms added and optimized by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, whereas Symm-Clashes lists symmetry related clashes.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	5549	0	5388	28	0
1	B	5538	0	5375	42	0
2	A	1	0	0	0	0
2	B	1	0	0	0	0
3	A	36	0	19	0	0
3	B	36	0	19	0	0
4	A	4	0	6	0	0
4	B	4	0	6	2	0
5	A	731	0	0	2	0
5	B	702	0	0	15	0
All	All	12602	0	10813	70	0

The all-atom clashscore is defined as the number of clashes found per 1000 atoms (including hydrogen atoms). The all-atom clashscore for this structure is 3.

All (70) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:212:MET:HE1	5:B:1588:HOH:O	1.58	1.02
4:B:805:EDO:O2	5:B:923:HOH:O	1.82	0.98
1:B:465:ASN:ND2	1:B:550:TYR:H	1.84	0.76
1:A:112:LEU:HD22	1:A:164:ILE:HB	1.70	0.74
1:B:465:ASN:HD21	1:B:550:TYR:H	1.40	0.69
1:A:636:GLN:HG2	5:B:1576:HOH:O	1.92	0.69
1:B:392:ARG:HH11	1:B:399:HIS:HE1	1.40	0.68
1:B:437:ALA:H	1:B:454:ASN:HD22	1.41	0.67
1:B:30:HIS:HD2	1:B:162:ASP:OD1	1.79	0.66
1:B:391:HIS:HD2	1:B:499:GLN:HE21	1.44	0.66
1:B:146:GLN:HE21	1:B:149:ASN:HD21	1.42	0.65
1:B:391:HIS:CD2	1:B:499:GLN:HE21	2.14	0.64
1:B:392:ARG:HH11	1:B:399:HIS:CE1	2.16	0.64
1:A:68:MET:HE3	1:A:111:TYR:HD1	1.64	0.63
1:B:391:HIS:HE1	1:B:501:SER:OG	1.81	0.63
1:A:128:SER:HB2	5:A:1599:HOH:O	2.00	0.61
1:A:392:ARG:HH11	1:A:399:HIS:HE1	1.48	0.60

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:89:HIS:HD2	1:B:122:TYR:OH	1.84	0.59
1:A:30:HIS:HD2	1:A:162:ASP:OD1	1.86	0.58
1:A:89:HIS:HD2	1:A:122:TYR:OH	1.86	0.58
1:A:392:ARG:HH11	1:A:399:HIS:CE1	2.22	0.57
1:B:201:ASN:HB2	5:B:1560:HOH:O	2.05	0.57
1:A:681:HIS:HE1	1:A:688:ASP:OD1	1.88	0.57
1:B:391:HIS:HD2	1:B:499:GLN:NE2	2.04	0.56
1:A:441:ASN:O	1:B:669:THR:HG21	2.06	0.55
1:A:252:TYR:H	1:A:259:GLN:NE2	2.04	0.55
1:B:70:GLU:HG3	5:B:1585:HOH:O	2.07	0.54
1:B:318:ILE:CG2	5:B:923:HOH:O	2.54	0.54
1:B:640:LEU:HB3	5:B:1599:HOH:O	2.07	0.54
1:B:252:TYR:H	1:B:259:GLN:NE2	2.06	0.52
1:A:392:ARG:HD3	1:A:399:HIS:CE1	2.44	0.52
1:A:612:SER:OG	1:A:681:HIS:HD2	1.93	0.51
1:B:396:GLY:O	1:B:399:HIS:HD2	1.94	0.51
1:B:389:SER:HB2	1:B:404:ALA:HB3	1.93	0.51
1:B:30:HIS:HE1	5:B:1128:HOH:O	1.94	0.50
1:B:350:SER:O	1:B:355:GLY:HA3	2.10	0.50
1:B:262:ALA:HA	5:B:1560:HOH:O	2.11	0.50
1:B:318:ILE:HG21	5:B:923:HOH:O	2.12	0.50
1:B:391:HIS:CD2	1:B:499:GLN:HB3	2.48	0.49
1:A:68:MET:HE3	1:A:111:TYR:CD1	2.47	0.49
1:B:451:LEU:HD13	1:B:633:ASN:HB3	1.95	0.49
1:B:78:LYS:HD2	5:B:1145:HOH:O	2.13	0.48
1:A:615:SER:HB2	1:A:647:ALA:HB2	1.96	0.48
1:B:201:ASN:HB3	1:B:252:TYR:OH	2.14	0.48
1:B:146:GLN:HE21	1:B:149:ASN:ND2	2.11	0.48
1:A:396:GLY:O	1:A:399:HIS:HD2	1.95	0.48
1:A:350:SER:O	1:A:355:GLY:HA3	2.15	0.47
1:B:624:GLU:HG2	1:B:638:GLN:HG2	1.96	0.46
1:A:409:GLN:O	1:A:414:GLY:HA3	2.15	0.46
1:B:304:GLY:HA2	5:B:923:HOH:O	2.15	0.46
1:A:389:SER:HB2	1:A:404:ALA:HB3	1.97	0.46
1:B:509:TYR:HB2	1:B:512:VAL:HB	1.99	0.45
1:B:318:ILE:HG23	5:B:923:HOH:O	2.15	0.45
1:A:90:LYS:HA	1:A:93:TYR:CE2	2.52	0.45
1:B:563:LYS:HE3	5:B:1090:HOH:O	2.17	0.44
1:A:259:GLN:HB2	1:A:259:GLN:HE21	1.69	0.43
1:B:90:LYS:HA	1:B:93:TYR:CE2	2.53	0.43
1:B:437:ALA:H	1:B:454:ASN:ND2	2.13	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:699:ASN:OD1	1:A:734:ARG:HD3	2.18	0.43
1:A:187:PHE:HA	1:A:191:GLU:HB2	2.01	0.43
1:A:509:TYR:HB2	1:A:512:VAL:HB	1.99	0.43
1:B:72:VAL:HG13	1:B:118:LEU:HD23	2.00	0.42
1:B:158:ILE:HG12	1:B:213:THR:HA	2.01	0.42
1:A:30:HIS:N	5:A:1613:HOH:O	2.51	0.42
1:A:418:LYS:O	1:A:419:LEU:HB2	2.19	0.42
1:B:318:ILE:HD13	4:B:805:EDO:H11	2.02	0.42
1:A:68:MET:CE	1:A:111:TYR:HD1	2.31	0.42
1:B:371:ASN:HB2	5:B:1265:HOH:O	2.19	0.42
1:A:391:HIS:O	1:A:401:ALA:HA	2.20	0.41
1:B:392:ARG:HD3	1:B:399:HIS:CE1	2.55	0.41

There are no symmetry-related clashes.

## 5.3 Torsion angles [i](#)

### 5.3.1 Protein backbone [i](#)

In the following table, the Percentiles column shows the percent Ramachandran outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

The Analysed column shows the number of residues for which the backbone conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	704/736 (96%)	682 (97%)	21 (3%)	1 (0%)	56	46
1	B	703/736 (96%)	679 (97%)	23 (3%)	1 (0%)	56	46
All	All	1407/1472 (96%)	1361 (97%)	44 (3%)	2 (0%)	56	46

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	718	ALA
1	B	719	GLY

### 5.3.2 Protein sidechains ⓘ

In the following table, the Percentiles column shows the percent sidechain outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

The Analysed column shows the number of residues for which the sidechain conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	586/610 (96%)	579 (99%)	7 (1%)	78	76
1	B	585/610 (96%)	576 (98%)	9 (2%)	72	69
All	All	1171/1220 (96%)	1155 (99%)	16 (1%)	74	71

All (16) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	112	LEU
1	A	135	ARG
1	A	150	GLU
1	A	499	GLN
1	A	535	LEU
1	A	656	LEU
1	A	724	GLU
1	B	56	GLN
1	B	135	ARG
1	B	143	LEU
1	B	150	GLU
1	B	311	ASP
1	B	499	GLN
1	B	640	LEU
1	B	656	LEU
1	B	720	GLU

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (27) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	30	HIS
1	A	89	HIS
1	A	174	GLN
1	A	259	GLN
1	A	399	HIS
1	A	428	ASN

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Mol	Chain	Res	Type
1	A	441	ASN
1	A	499	GLN
1	A	619	ASN
1	A	636	GLN
1	A	681	HIS
1	B	30	HIS
1	B	56	GLN
1	B	89	HIS
1	B	149	ASN
1	B	259	GLN
1	B	337	ASN
1	B	344	GLN
1	B	391	HIS
1	B	399	HIS
1	B	413	HIS
1	B	454	ASN
1	B	460	GLN
1	B	465	ASN
1	B	499	GLN
1	B	636	GLN
1	B	675	ASN

### 5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

### 5.4 Non-standard residues in protein, DNA, RNA chains ⓘ

There are no non-standard protein/DNA/RNA residues in this entry.

### 5.5 Carbohydrates ⓘ

6 carbohydrates are modelled in this entry.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the chemical component dictionary. The Link column lists molecule types, if any, to which the group is linked. The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with  $|Z| > 2$  is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z  > 2	Counts	RMSZ	# Z  > 2
3	LGU	A	802	3	10,13,13	0.57	0	15,19,19	0.91	2 (13%)
3	MAV	A	803	3	9,12,13	0.52	0	13,17,19	0.95	1 (7%)
3	BEM	A	804	3	8,11,13	0.57	0	8,15,19	1.34	1 (12%)
3	LGU	B	802	3	10,13,13	0.57	0	15,19,19	0.97	0
3	MAV	B	803	3	9,12,13	0.53	0	13,17,19	0.70	0
3	BEM	B	804	3	8,11,13	0.43	0	8,15,19	0.97	1 (12%)

In the following table, the Chirals column lists the number of chiral outliers, the number of chiral centers analysed, the number of these observed in the model and the number defined in the chemical component dictionary. Similar counts are reported in the Torsion and Rings columns. '-' means no outliers of that kind were identified.

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
3	LGU	A	802	3	-	0/0/24/24	0/1/1/1
3	MAV	A	803	3	1/1/5/6	0/0/21/24	0/1/1/1
3	BEM	A	804	3	-	0/0/17/24	0/1/1/1
3	LGU	B	802	3	-	0/0/24/24	0/1/1/1
3	MAV	B	803	3	1/1/5/6	0/0/21/24	0/1/1/1
3	BEM	B	804	3	-	0/0/17/24	0/1/1/1

There are no bond length outliers.

All (5) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	803	MAV	O4-C4-C5	-2.29	106.10	110.42
3	A	802	LGU	C1-C2-C3	2.10	113.55	110.43
3	A	802	LGU	C1-O5-C5	2.20	115.46	112.22
3	B	804	BEM	C4-C3-C2	2.30	113.28	110.23
3	A	804	BEM	C4-C3-C2	3.09	114.33	110.23

All (2) chirality outliers are listed below:

Mol	Chain	Res	Type	Atom
3	B	803	MAV	C1
3	A	803	MAV	C1

There are no torsion outliers.

There are no ring outliers.

No monomer is involved in short contacts.

## 5.6 Ligand geometry

Of 4 ligands modelled in this entry, 2 are monoatomic - leaving 2 for Mogul analysis.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the chemical component dictionary. The Link column lists molecule types, if any, to which the group is linked. The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with  $|Z| > 2$  is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\# Z  > 2$	Counts	RMSZ	$\# Z  > 2$
4	EDO	A	805	-	3,3,3	0.47	0	2,2,2	0.41	0
4	EDO	B	805	-	3,3,3	0.40	0	2,2,2	0.63	0

In the following table, the Chirals column lists the number of chiral outliers, the number of chiral centers analysed, the number of these observed in the model and the number defined in the chemical component dictionary. Similar counts are reported in the Torsion and Rings columns. '-' means no outliers of that kind were identified.

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
4	EDO	A	805	-	-	0/1/1/1	0/0/0/0
4	EDO	B	805	-	-	0/1/1/1	0/0/0/0

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

1 monomer is involved in 2 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	B	805	EDO	2	0

## 5.7 Other polymers

There are no such residues in this entry.

## 5.8 Polymer linkage issues ⓘ

There are no chain breaks in this entry.



## 6 Fit of model and data [i](#)

### 6.1 Protein, DNA and RNA chains [i](#)

In the following table, the column labelled ‘#RSRZ> 2’ contains the number (and percentage) of RSRZ outliers, followed by percent RSRZ outliers for the chain as percentile scores relative to all X-ray entries and entries of similar resolution. The OWAB column contains the minimum, median, 95<sup>th</sup> percentile and maximum values of the occupancy-weighted average B-factor per residue. The column labelled ‘Q< 0.9’ lists the number of (and percentage) of residues with an average occupancy less than 0.9.

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å <sup>2</sup> )	Q<0.9
1	A	706/736 (95%)	-0.29	7 (0%) 84 86	12, 19, 32, 77	0
1	B	705/736 (95%)	-0.16	12 (1%) 73 76	13, 20, 33, 75	0
All	All	1411/1472 (95%)	-0.23	19 (1%) 79 82	12, 19, 33, 77	0

All (19) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	720	GLU	6.6
1	B	719	GLY	5.2
1	B	720	GLU	4.4
1	B	721	GLU	4.4
1	A	719	GLY	4.1
1	B	718	ALA	3.9
1	A	718	ALA	3.0
1	B	51	ASP	2.9
1	B	722	ALA	2.9
1	B	713	VAL	2.8
1	A	724	GLU	2.5
1	B	717	ASN	2.5
1	B	138	GLN	2.4
1	A	721	GLU	2.4
1	A	51	ASP	2.3
1	B	715	LYS	2.3
1	B	363	GLY	2.3
1	B	724	GLU	2.1
1	A	138	GLN	2.1

### 6.2 Non-standard residues in protein, DNA, RNA chains [i](#)

There are no non-standard protein/DNA/RNA residues in this entry.

### 6.3 Carbohydrates [i](#)

In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. LLDF column lists the quality of electron density of the group with respect to its neighbouring residues in protein, DNA or RNA chains. The B-factors column lists the minimum, median, 95<sup>th</sup> percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors(Å <sup>2</sup> )	Q<0.9
3	BEM	A	804	11/13	0.96	0.12	1.03	19,20,21,21	0
3	MAV	A	803	12/13	0.92	0.11	-0.01	22,25,29,30	0
3	BEM	B	804	11/13	0.95	0.11	-0.06	18,19,20,20	0
3	LGU	A	802	13/13	0.92	0.10	-0.35	28,30,33,34	0
3	MAV	B	803	12/13	0.94	0.09	-0.42	21,24,29,30	0
3	LGU	B	802	13/13	0.92	0.10	-0.46	27,30,34,35	0

### 6.4 Ligands [i](#)

In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. LLDF column lists the quality of electron density of the group with respect to its neighbouring residues in protein, DNA or RNA chains. The B-factors column lists the minimum, median, 95<sup>th</sup> percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors(Å <sup>2</sup> )	Q<0.9
4	EDO	B	805	4/4	0.84	0.20	4.75	24,25,27,28	0
4	EDO	A	805	4/4	0.97	0.07	-0.83	16,17,17,17	0
2	ZN	A	801	1/1	1.00	0.05	-1.81	17,17,17,17	0
2	ZN	B	801	1/1	1.00	0.05	-2.08	15,15,15,15	0

### 6.5 Other polymers [i](#)

There are no such residues in this entry.